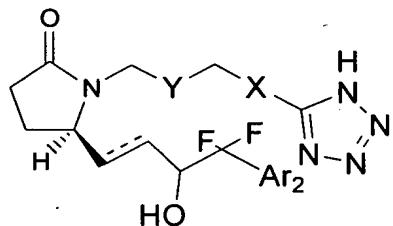


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In the Claims

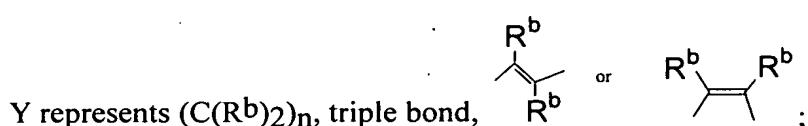
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17. A compound of structural formula I:



I

or a pharmaceutically acceptable salt, enantiomer, diastereomer, pro drug or mixture thereof, wherein

X is $(CH_2)_n$, O or S;



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Ar₂ independently represent (CH₂)_mC₆₋₁₀aryl, (CH₂)_mC₅₋₁₀heteroaryl, (CH₂)_mC₃₋₁₀ heterocycloalkyl, (CH₂)_mC₃₋₈ cycloalkyl said cycloalkyl, heterocycloalkyl, aryl or heteroaryl unsubstituted or substituted with 1-3 groups of R_a;

R_a represents C₁₋₆ alkoxy, C₁₋₆ alkyl, CF₃, nitro, amino, cyano, C₁₋₆ alkylamino, or halogen;

R^b independently represents H, halogen, C₁₋₆ alkyl, C₃₋₆ cylcoalkyl or

— represents a double or single bond;

n represents 0-4; and

m represents 0-8.

18. The compound according to claim 17 wherein X and Y are (CH₂)_n, — represents a double bond; and Ar₂ is phenyl.

19. The compound according to claim 18 wherein X is (CH₂)_n and n is 1 and Y is (CH₂)_n and n is 3.